



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 08:59 AM EDT

PDB ID : 4DUP / pdb_00004dup
Title : Crystal Structure of a quinone oxidoreductase from Rhizobium etli CFN 42
Authors : Kumaran, D.; Rice, S.; Evans, B.; Foti, R.; Gizzi, A.; Hillerich, B.; Kar, A.; Lafleur, J.; Seidel, R.; Villigas, G.; Zencheck, W.; Almo, S.C.; Swaminathan, S.; New York Structural Genomics Research Consortium (NYSGRG)
Deposited on : 2012-02-22
Resolution : 2.45 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

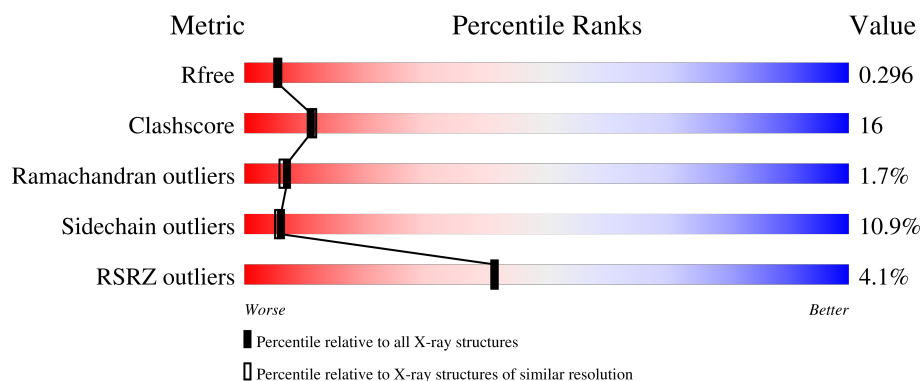
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	353	 % 65% 25% 6%
1	B	353	 6% 61% 23% 12%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 4821 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called quinone oxidoreductase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	Se	0	0	0
			2442	1553	417	459	5	8			
1	B	312	Total	C	N	O	S	Se	0	0	0
			2322	1485	395	431	4	7			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MSE	-	expression tag	UNP Q2KB38
A	-20	HIS	-	expression tag	UNP Q2KB38
A	-19	HIS	-	expression tag	UNP Q2KB38
A	-18	HIS	-	expression tag	UNP Q2KB38
A	-17	HIS	-	expression tag	UNP Q2KB38
A	-16	HIS	-	expression tag	UNP Q2KB38
A	-15	HIS	-	expression tag	UNP Q2KB38
A	-14	SER	-	expression tag	UNP Q2KB38
A	-13	SER	-	expression tag	UNP Q2KB38
A	-12	GLY	-	expression tag	UNP Q2KB38
A	-11	VAL	-	expression tag	UNP Q2KB38
A	-10	ASP	-	expression tag	UNP Q2KB38
A	-9	LEU	-	expression tag	UNP Q2KB38
A	-8	GLY	-	expression tag	UNP Q2KB38
A	-7	THR	-	expression tag	UNP Q2KB38
A	-6	GLU	-	expression tag	UNP Q2KB38
A	-5	ASN	-	expression tag	UNP Q2KB38
A	-4	LEU	-	expression tag	UNP Q2KB38
A	-3	TYR	-	expression tag	UNP Q2KB38
A	-2	PHE	-	expression tag	UNP Q2KB38
A	-1	GLN	-	expression tag	UNP Q2KB38
A	0	SER	-	expression tag	UNP Q2KB38
B	-21	MSE	-	expression tag	UNP Q2KB38
B	-20	HIS	-	expression tag	UNP Q2KB38
B	-19	HIS	-	expression tag	UNP Q2KB38

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	HIS	-	expression tag	UNP Q2KB38
B	-17	HIS	-	expression tag	UNP Q2KB38
B	-16	HIS	-	expression tag	UNP Q2KB38
B	-15	HIS	-	expression tag	UNP Q2KB38
B	-14	SER	-	expression tag	UNP Q2KB38
B	-13	SER	-	expression tag	UNP Q2KB38
B	-12	GLY	-	expression tag	UNP Q2KB38
B	-11	VAL	-	expression tag	UNP Q2KB38
B	-10	ASP	-	expression tag	UNP Q2KB38
B	-9	LEU	-	expression tag	UNP Q2KB38
B	-8	GLY	-	expression tag	UNP Q2KB38
B	-7	THR	-	expression tag	UNP Q2KB38
B	-6	GLU	-	expression tag	UNP Q2KB38
B	-5	ASN	-	expression tag	UNP Q2KB38
B	-4	LEU	-	expression tag	UNP Q2KB38
B	-3	TYR	-	expression tag	UNP Q2KB38
B	-2	PHE	-	expression tag	UNP Q2KB38
B	-1	GLN	-	expression tag	UNP Q2KB38
B	0	SER	-	expression tag	UNP Q2KB38

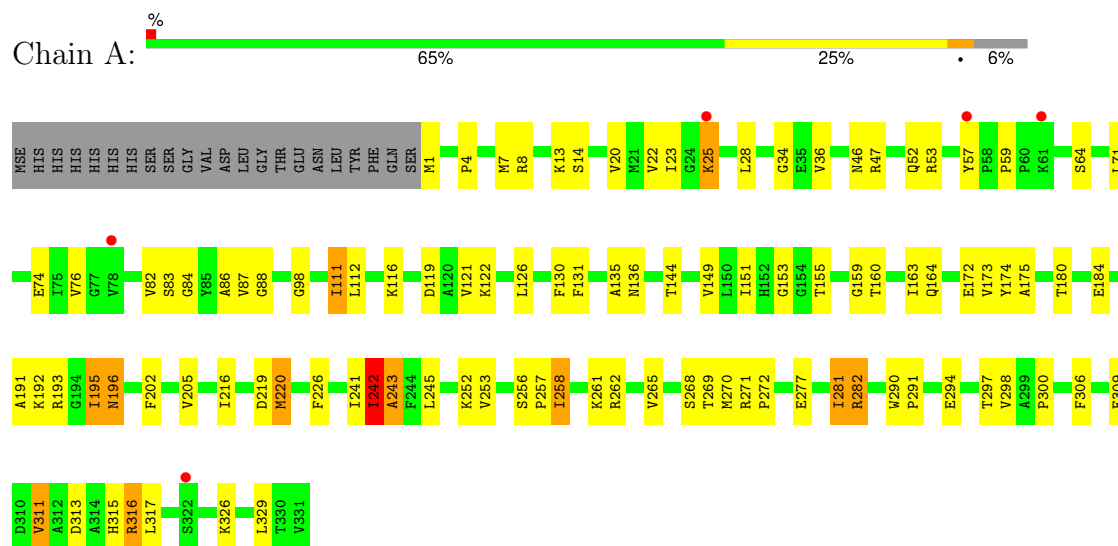
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	30	Total O 30 30	0	0
2	B	27	Total O 27 27	0	0

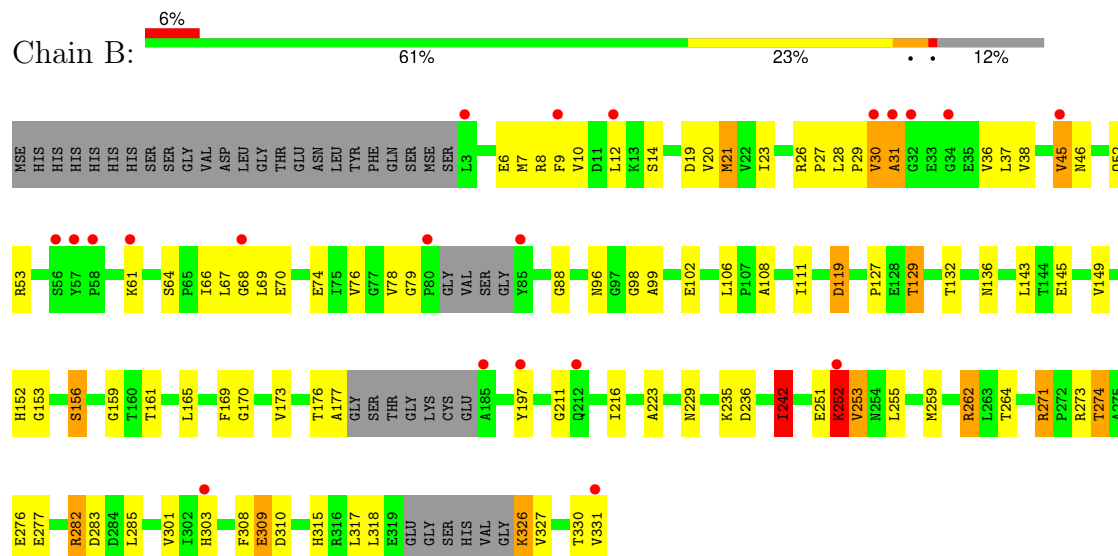
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: quinone oxidoreductase



- Molecule 1: quinone oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.71Å 142.82Å 59.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.36 – 2.45 47.36 – 2.45	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.36-2.45) 99.6 (47.36-2.45)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.52 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.222 , 0.293 0.226 , 0.296	Depositor DCC
R_{free} test set	1495 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	45.9	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4821	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.04	2/2480 (0.1%)	1.26	16/3350 (0.5%)
1	B	0.99	3/2356 (0.1%)	1.24	7/3181 (0.2%)
All	All	1.02	5/4836 (0.1%)	1.25	23/6531 (0.4%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	303	HIS	CG-CD2	6.13	1.42	1.35
1	A	47	ARG	C-O	-5.78	1.19	1.24
1	B	303	HIS	CB-CG	5.39	1.57	1.50
1	A	243	ALA	N-CA	5.35	1.52	1.46
1	B	315	HIS	CG-CD2	5.15	1.41	1.35

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	173	VAL	N-CA-C	9.52	121.43	108.11
1	B	242	ILE	CB-CA-C	9.17	119.64	110.65
1	A	311	VAL	CB-CA-C	-8.89	100.50	111.88
1	A	242	ILE	N-CA-C	-7.97	105.72	113.53
1	B	21	MSE	CG-SE-CE	-7.75	81.88	98.92
1	B	253	VAL	N-CA-C	7.12	118.14	107.75
1	A	111	ILE	CB-CA-C	-6.99	101.42	111.34
1	A	270	MSE	N-CA-C	6.79	118.37	110.97
1	A	242	ILE	N-CA-CB	-6.54	103.29	112.12
1	A	59	PRO	O-C-N	6.43	124.27	121.31
1	A	281	ILE	N-CA-C	-6.41	104.25	110.72
1	A	13	LYS	N-CA-C	-5.94	104.72	111.07
1	A	220	MSE	CB-CA-C	-5.58	99.95	109.65
1	B	262	ARG	CB-CA-C	-5.42	104.19	111.89
1	B	242	ILE	N-CA-CB	-5.33	104.93	112.12
1	A	98	GLY	N-CA-C	5.31	121.01	114.69
1	A	131	PHE	N-CA-C	-5.31	105.39	111.07

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	223	ALA	N-CA-C	5.30	117.48	111.02
1	A	84	GLY	N-CA-C	5.29	122.91	115.30
1	A	242	ILE	CG1-CB-CG2	-5.15	95.25	110.70
1	A	173	VAL	N-CA-C	5.10	115.83	108.48
1	A	265	VAL	N-CA-C	-5.09	101.05	108.17
1	A	311	VAL	N-CA-CB	5.05	116.11	110.51

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2442	0	2478	69	0
1	B	2322	0	2363	103	0
2	A	30	0	0	0	0
2	B	27	0	0	1	0
All	All	4821	0	4841	158	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 16.

All (158) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:MSE:CE	1:B:29:PRO:HD2	1.58	1.34
1:B:45:VAL:HG22	1:B:327:VAL:HG23	1.28	1.16
1:B:271:ARG:HH11	1:B:271:ARG:HG3	1.13	1.10
1:B:282:ARG:HG2	1:B:282:ARG:HH11	1.05	1.09
1:A:282:ARG:HH11	1:A:282:ARG:HG2	1.01	1.08
1:B:7:MSE:HE1	1:B:29:PRO:HD2	1.12	1.08
1:B:7:MSE:CE	1:B:29:PRO:CD	2.37	1.02
1:A:243:ALA:N	1:B:259:MSE:HE1	1.76	0.99
1:B:7:MSE:CE	1:B:28:LEU:HA	1.93	0.98
1:B:7:MSE:HE1	1:B:29:PRO:CD	1.93	0.96

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:ARG:HG2	1:B:282:ARG:NH1	1.78	0.96
1:B:7:MSE:HE3	1:B:28:LEU:HA	1.46	0.96
1:B:111:ILE:O	1:B:282:ARG:HD2	1.64	0.95
1:A:111:ILE:O	1:A:282:ARG:HD2	1.66	0.95
1:A:53:ARG:HH11	1:A:315:HIS:HD2	0.99	0.94
1:A:53:ARG:HH11	1:A:315:HIS:CD2	1.87	0.93
1:A:241:ILE:HG22	1:B:259:MSE:HE3	1.48	0.92
1:A:282:ARG:HG2	1:A:282:ARG:NH1	1.80	0.92
1:A:242:ILE:C	1:B:259:MSE:CE	2.43	0.91
1:B:274:THR:HG22	1:B:277:GLU:H	1.34	0.91
1:B:21:MSE:CE	1:B:67:LEU:CD1	2.49	0.91
1:B:45:VAL:CG2	1:B:327:VAL:HG23	2.03	0.88
1:A:242:ILE:C	1:B:259:MSE:HE2	2.00	0.85
1:A:149:VAL:HG12	1:A:216:ILE:HB	1.58	0.84
1:B:271:ARG:HH11	1:B:271:ARG:CG	1.89	0.84
1:B:271:ARG:HG3	1:B:271:ARG:NH1	1.83	0.84
1:B:282:ARG:HH11	1:B:282:ARG:CG	1.88	0.83
1:A:282:ARG:HH11	1:A:282:ARG:CG	1.87	0.82
1:A:243:ALA:CA	1:B:259:MSE:HE1	2.10	0.81
1:B:274:THR:HG22	1:B:277:GLU:N	1.97	0.80
1:B:153:GLY:O	1:B:156:SER:HB2	1.80	0.80
1:A:135:ALA:HB2	1:A:281:ILE:HD12	1.62	0.79
1:B:7:MSE:HE3	1:B:29:PRO:CD	2.12	0.78
1:A:242:ILE:C	1:B:259:MSE:HE1	2.06	0.78
1:A:53:ARG:NH1	1:A:315:HIS:HD2	1.77	0.78
1:B:7:MSE:HE2	1:B:28:LEU:HG	1.67	0.76
1:B:7:MSE:HE3	1:B:29:PRO:HD2	1.66	0.76
1:A:242:ILE:CA	1:B:259:MSE:HE2	2.16	0.75
1:B:7:MSE:HE2	1:B:28:LEU:HA	1.70	0.73
1:B:66:ILE:HB	1:B:98:GLY:HA2	1.69	0.73
1:B:8:ARG:NH2	1:B:309:GLU:OE2	2.22	0.72
1:A:243:ALA:N	1:B:259:MSE:CE	2.51	0.71
1:A:52:GLN:NE2	1:A:271:ARG:HH12	1.88	0.70
1:B:156:SER:HB3	1:B:159:GLY:H	1.57	0.70
1:B:26:ARG:HB3	1:B:27:PRO:HD2	1.72	0.69
1:B:111:ILE:O	1:B:282:ARG:CD	2.39	0.69
1:B:21:MSE:HE1	1:B:67:LEU:HD13	1.73	0.69
1:A:126:LEU:HD23	1:A:300:PRO:HG3	1.75	0.69
1:B:317:LEU:HD23	1:B:327:VAL:HG11	1.76	0.68
1:B:21:MSE:HE1	1:B:67:LEU:CD1	2.22	0.68
1:A:52:GLN:HE22	1:A:271:ARG:HH12	1.43	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:ASP:OD2	1:A:121:VAL:HG22	1.95	0.67
1:B:236:ASP:OD1	1:B:262:ARG:HD2	1.95	0.67
1:B:30:VAL:HG13	1:B:31:ALA:H	1.60	0.65
1:B:21:MSE:HE3	1:B:67:LEU:CD1	2.24	0.65
1:B:70:GLU:HG2	1:B:127:PRO:HD2	1.79	0.64
1:B:64:SER:OG	1:B:66:ILE:HD12	1.97	0.64
1:B:152:HIS:HE1	1:B:229:ASN:OD1	1.83	0.62
1:A:243:ALA:HA	1:B:259:MSE:HE1	1.81	0.62
1:A:126:LEU:CD2	1:A:300:PRO:HG3	2.30	0.62
1:A:7:MSE:HB2	1:A:28:LEU:HD12	1.81	0.62
1:B:68:GLY:HA2	1:B:99:ALA:HB3	1.83	0.60
1:A:112:LEU:HD13	1:A:290:TRP:CH2	2.36	0.60
1:B:282:ARG:NH1	1:B:282:ARG:CG	2.54	0.60
1:A:172:GLU:HG3	1:A:192:LYS:HD3	1.83	0.59
1:B:78:VAL:HG12	1:B:79:GLY:N	2.17	0.59
1:B:21:MSE:CE	1:B:67:LEU:HD12	2.31	0.59
1:A:277:GLU:O	1:A:281:ILE:HG12	2.04	0.58
1:A:253:VAL:HG23	1:B:253:VAL:CG1	2.34	0.57
1:A:53:ARG:NH1	1:A:315:HIS:CD2	2.60	0.57
1:B:12:LEU:HD23	1:B:20:VAL:HG13	1.86	0.57
1:A:253:VAL:CG2	1:B:253:VAL:HG11	2.34	0.57
1:B:21:MSE:HE3	1:B:67:LEU:HD11	1.85	0.57
1:B:8:ARG:HE	1:B:23:ILE:HG21	1.69	0.57
1:B:129:THR:CG2	1:B:161:THR:OG1	2.54	0.56
1:B:21:MSE:CE	1:B:67:LEU:HD13	2.30	0.55
1:B:143:LEU:HD13	1:B:216:ILE:HD12	1.89	0.55
1:B:326:LYS:N	2:B:415:HOH:O	2.39	0.55
1:B:30:VAL:O	1:B:31:ALA:CB	2.54	0.55
1:A:226:PHE:HE2	1:A:253:VAL:HG13	1.73	0.53
1:A:253:VAL:HG21	1:B:253:VAL:HG11	1.90	0.53
1:B:7:MSE:HE3	1:B:29:PRO:HD3	1.90	0.53
1:A:111:ILE:O	1:A:282:ARG:CD	2.49	0.53
1:B:145:GLU:HG3	1:B:170:GLY:HA3	1.89	0.53
1:A:294:GLU:HA	1:A:294:GLU:OE1	2.08	0.53
1:A:253:VAL:CG2	1:B:253:VAL:CG1	2.87	0.52
1:A:195:ILE:N	1:A:195:ILE:HD13	2.25	0.52
1:A:46:ASN:HD21	1:A:326:LYS:NZ	2.08	0.52
1:B:108:ALA:HA	1:B:111:ILE:HD12	1.92	0.52
1:B:282:ARG:NH1	1:B:283:ASP:OD1	2.43	0.52
1:A:136:ASN:ND2	1:A:268:SER:OG	2.42	0.51
1:B:274:THR:HG23	1:B:276:GLU:H	1.76	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:VAL:CG2	1:B:327:VAL:CG2	2.86	0.50
1:B:26:ARG:CB	1:B:27:PRO:HD2	2.38	0.50
1:B:301:VAL:O	1:B:326:LYS:HB2	2.12	0.50
1:A:271:ARG:HB2	1:A:272:PRO:HD3	1.94	0.50
1:A:34:GLY:HA2	1:A:82:VAL:HG22	1.93	0.49
1:A:258:ILE:N	1:A:258:ILE:HD13	2.27	0.49
1:B:68:GLY:HA2	1:B:99:ALA:H	1.78	0.49
1:B:177:ALA:O	1:B:197:TYR:HD2	1.96	0.49
1:A:261:LYS:O	1:A:262:ARG:HB2	2.13	0.49
1:B:273:ARG:HG2	1:B:277:GLU:OE1	2.13	0.48
1:B:274:THR:HG23	1:B:276:GLU:N	2.27	0.48
1:A:282:ARG:NH1	1:A:282:ARG:CG	2.52	0.48
1:B:12:LEU:HD21	1:B:53:ARG:HA	1.94	0.48
1:B:52:GLN:HE22	1:B:271:ARG:HH22	1.60	0.48
1:B:78:VAL:HG12	1:B:79:GLY:H	1.78	0.48
1:A:7:MSE:O	1:A:25:LYS:HA	2.14	0.48
1:A:195:ILE:HD12	1:A:205:VAL:HG11	1.96	0.48
1:A:74:GLU:OE2	1:A:88:GLY:HA2	2.14	0.47
1:B:26:ARG:HB3	1:B:27:PRO:CD	2.42	0.47
1:A:220:MSE:HA	1:A:242:ILE:HG22	1.97	0.47
1:A:313:ASP:OD1	1:A:316:ARG:NH1	2.48	0.47
1:A:153:GLY:O	1:A:159:GLY:HA3	2.14	0.47
1:A:195:ILE:HG21	1:A:202:PHE:HA	1.97	0.47
1:A:196:ASN:OD1	1:A:196:ASN:C	2.56	0.46
1:A:226:PHE:CE2	1:A:253:VAL:HG13	2.50	0.46
1:B:7:MSE:CE	1:B:29:PRO:HD3	2.41	0.46
1:B:74:GLU:OE2	1:B:88:GLY:HA2	2.16	0.46
1:A:151:ILE:O	1:A:175:ALA:HA	2.16	0.46
1:B:12:LEU:HD23	1:B:20:VAL:CG1	2.45	0.46
1:A:14:SER:O	1:A:20:VAL:HG21	2.15	0.46
1:A:7:MSE:HG2	1:A:8:ARG:O	2.15	0.46
1:A:291:PRO:O	1:A:294:GLU:HB2	2.16	0.45
1:B:45:VAL:HG22	1:B:327:VAL:CG2	2.20	0.45
1:B:274:THR:CG2	1:B:277:GLU:H	2.16	0.45
1:A:86:ALA:O	1:A:87:VAL:C	2.59	0.45
1:A:219:ASP:O	1:A:242:ILE:HG22	2.16	0.45
1:B:8:ARG:NH1	1:B:102:GLU:OE2	2.50	0.45
1:B:8:ARG:NH2	1:B:308:PHE:HD2	2.15	0.45
1:B:177:ALA:O	1:B:197:TYR:CD2	2.70	0.45
1:B:251:GLU:O	1:B:252:LYS:C	2.59	0.45
1:A:22:VAL:HG22	1:A:23:ILE:N	2.33	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:271:ARG:CG	1:B:271:ARG:NH1	2.59	0.44
1:B:26:ARG:CB	1:B:27:PRO:CD	2.97	0.43
1:B:10:VAL:HG11	1:B:21:MSE:HE2	2.01	0.42
1:A:242:ILE:HG21	1:A:242:ILE:HD13	1.14	0.42
1:A:256:SER:HB3	1:A:257:PRO:HD3	2.01	0.42
1:B:9:PHE:CE1	1:B:26:ARG:HB2	2.55	0.42
1:B:165:LEU:O	1:B:169:PHE:HD1	2.03	0.42
1:A:306:PHE:O	1:A:329:LEU:HA	2.18	0.42
1:A:191:ALA:O	1:A:192:LYS:C	2.63	0.42
1:A:180:THR:HG22	1:A:184:GLU:OE2	2.20	0.42
1:A:192:LYS:HE2	1:A:192:LYS:HB3	1.86	0.41
1:A:294:GLU:OE1	1:A:294:GLU:CA	2.68	0.41
1:B:285:LEU:HD23	1:B:285:LEU:HA	1.90	0.41
1:B:46:ASN:HD21	1:B:326:LYS:NZ	2.17	0.41
1:B:38:VAL:HG21	1:B:106:LEU:HD12	2.03	0.41
1:B:136:ASN:CG	1:B:242:ILE:HD11	2.44	0.41
1:A:126:LEU:O	1:A:130:PHE:HB2	2.20	0.41
1:A:172:GLU:OE2	1:A:174:TYR:OH	2.33	0.41
1:A:242:ILE:N	1:B:259:MSE:CE	2.84	0.41
1:B:149:VAL:O	1:B:149:VAL:HG23	2.21	0.41
1:B:78:VAL:CG1	1:B:79:GLY:N	2.82	0.40
1:B:136:ASN:ND2	1:B:242:ILE:HD11	2.37	0.40
1:B:7:MSE:HE2	1:B:28:LEU:CA	2.46	0.40
1:B:331:VAL:HG23	1:B:331:VAL:O	2.21	0.40
1:B:129:THR:HG21	1:B:161:THR:OG1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	329/353 (93%)	316 (96%)	12 (4%)	1 (0%)	36	45

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	304/353 (86%)	278 (91%)	16 (5%)	10 (3%)	3	1
All	All	633/706 (90%)	594 (94%)	28 (4%)	11 (2%)	7	6

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	14	SER
1	B	31	ALA
1	B	119	ASP
1	B	69	LEU
1	B	96	ASN
1	B	211	GLY
1	B	252	LYS
1	B	318	LEU
1	B	310	ASP
1	B	30	VAL
1	A	4	PRO

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	250/261 (96%)	220 (88%)	30 (12%)	5	4
1	B	237/261 (91%)	214 (90%)	23 (10%)	8	8
All	All	487/522 (93%)	434 (89%)	53 (11%)	6	5

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	25	LYS
1	A	36	VAL
1	A	57	TYR
1	A	64	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	71	LEU
1	A	76	VAL
1	A	83	SER
1	A	116	LYS
1	A	122	LYS
1	A	144	THR
1	A	155	THR
1	A	160	THR
1	A	163	ILE
1	A	164	GLN
1	A	193	ARG
1	A	195	ILE
1	A	196	ASN
1	A	242	ILE
1	A	245	LEU
1	A	252	LYS
1	A	258	ILE
1	A	269	THR
1	A	282	ARG
1	A	297	THR
1	A	298	VAL
1	A	309	GLU
1	A	311	VAL
1	A	316	ARG
1	A	317	LEU
1	B	6	GLU
1	B	19	ASP
1	B	36	VAL
1	B	37	LEU
1	B	45	VAL
1	B	61	LYS
1	B	76	VAL
1	B	119	ASP
1	B	129	THR
1	B	132	THR
1	B	156	SER
1	B	176	THR
1	B	235	LYS
1	B	242	ILE
1	B	252	LYS
1	B	255	LEU
1	B	264	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	271	ARG
1	B	274	THR
1	B	282	ARG
1	B	309	GLU
1	B	326	LYS
1	B	330	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	52	GLN
1	A	136	ASN
1	A	315	HIS
1	B	46	ASN
1	B	52	GLN
1	B	96	ASN
1	B	152	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	323/353 (91%)	0.14	5 (1%) 72 74	23, 43, 66, 84	0
1	B	305/353 (86%)	0.35	21 (6%) 23 20	22, 47, 74, 105	0
All	All	628/706 (88%)	0.24	26 (4%) 41 41	22, 45, 70, 105	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	57	TYR	4.3
1	B	185	ALA	3.5
1	B	56	SER	3.5
1	B	80	PRO	3.4
1	B	61	LYS	3.2
1	B	68	GLY	3.1
1	B	30	VAL	2.9
1	B	303	HIS	2.8
1	B	85	TYR	2.8
1	A	57	TYR	2.7
1	B	34	GLY	2.5
1	A	61	LYS	2.5
1	B	331	VAL	2.3
1	B	58	PRO	2.2
1	B	32	GLY	2.2
1	B	12	LEU	2.2
1	A	322	SER	2.2
1	B	212	GLN	2.1
1	B	45	VAL	2.1
1	A	25	LYS	2.1
1	B	9	PHE	2.1
1	B	252	LYS	2.1
1	B	197	TYR	2.1
1	A	78	VAL	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	3	LEU	2.0
1	B	31	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.